

## **Interzeolite conversion of micronsized FAU to nanosized CHA zeolite free of organic structure directing agent with high a CO<sub>2</sub> capacity**

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## Supporting information

**Table S1: List of conditions for selected synthesis samples**

Entry #	Name	FAU source	H <sub>2</sub> O/Si	K/Si	Na/Si	Time / h	Phase	Yield <sup>a</sup>
1	CHA(1.9)	CBV400	40	0.76	0.25	96	CHA*	113%
2		CBV400	40	0.51	0.51	96	CHA + minor FAU	
3		CBV720	30	0.5	-	168	Amorphous	
4	CHA(2.3)	CBV720	30	1	-	168	CHA**	24%

<sup>a</sup>Yield calculated by (mass product)/(mass starting zeolite) \*Minor impurity of LTA or GME

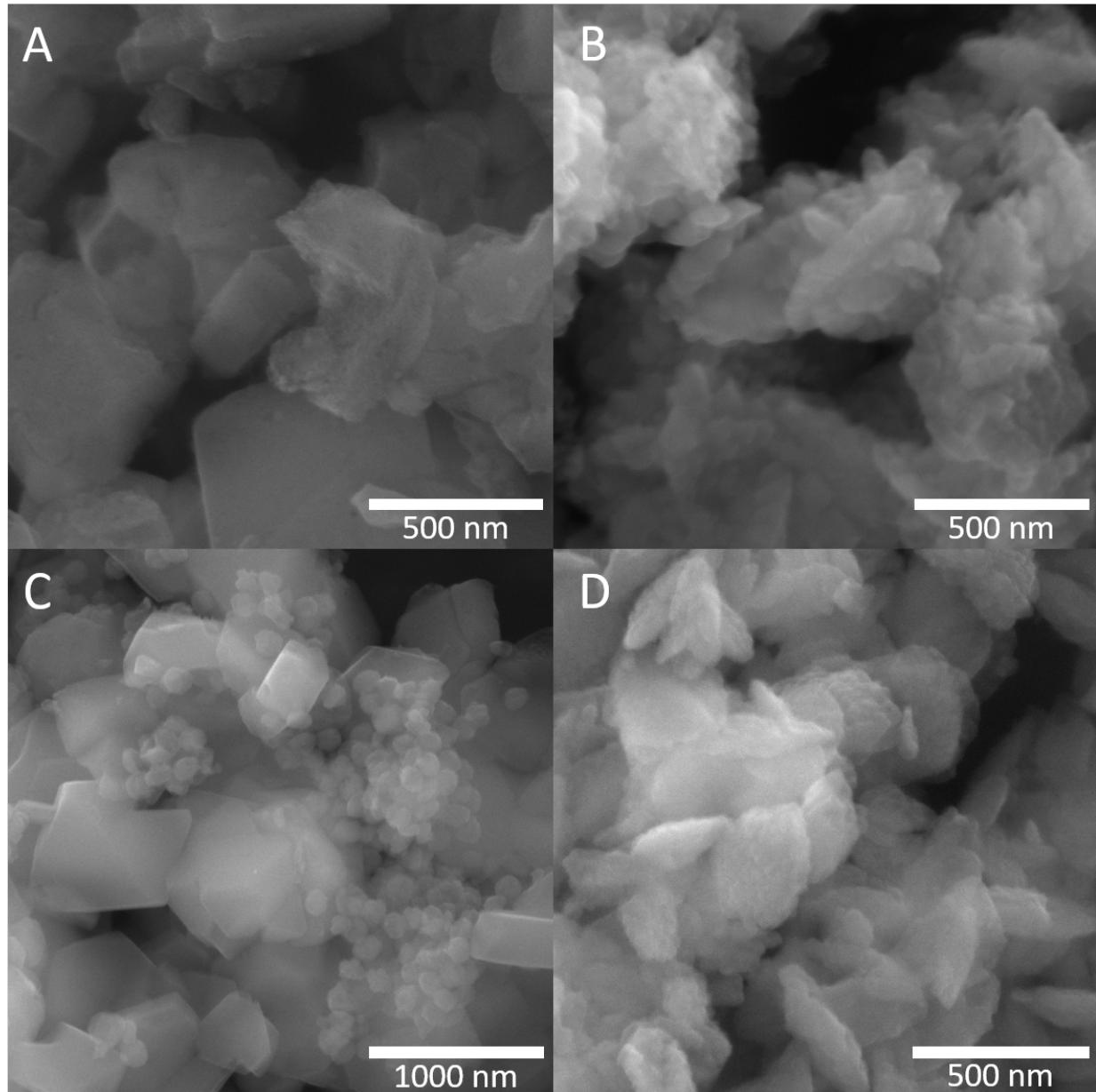
\*\*Minor impurity of LTL

### ICP analysis

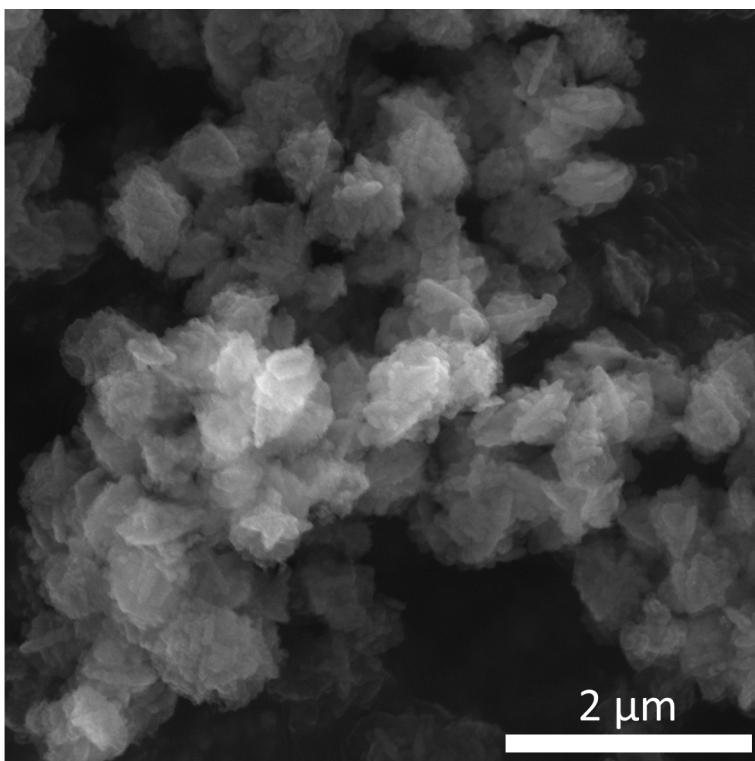
**Table S2: ICP results for CHA zeolite samples**

Sample	Si/Al	Na/Al	K/Al
CHA(1.9)	1.9	0.06	0.93
CHA(2.3)	2.3	0.01	0.97

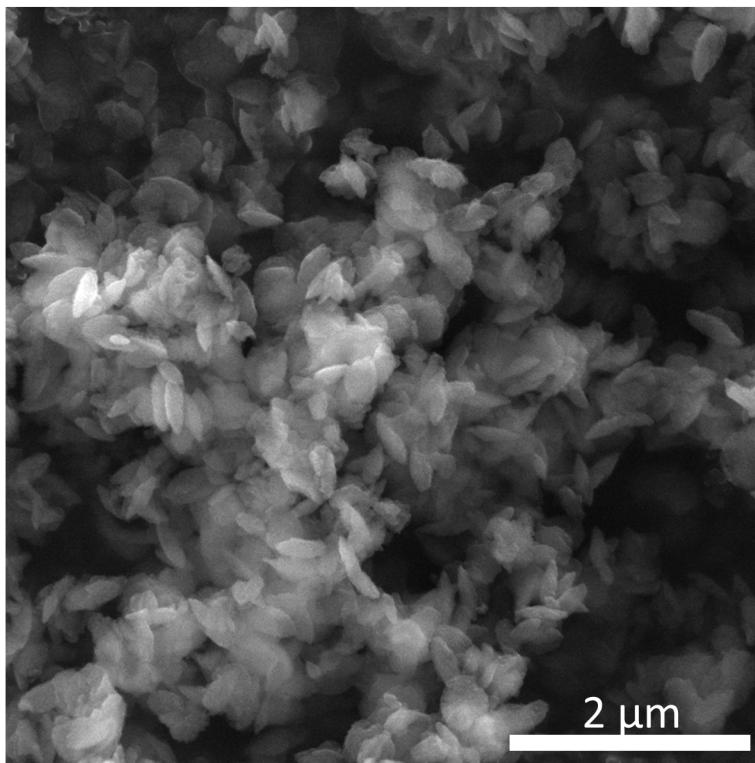
**SEM images**



**Figure S1: SEM images of (A) FAU(15), (B) CHA(2.3), (C) FAU(2.6), and (D) CHA(1.9) zeolite samples**

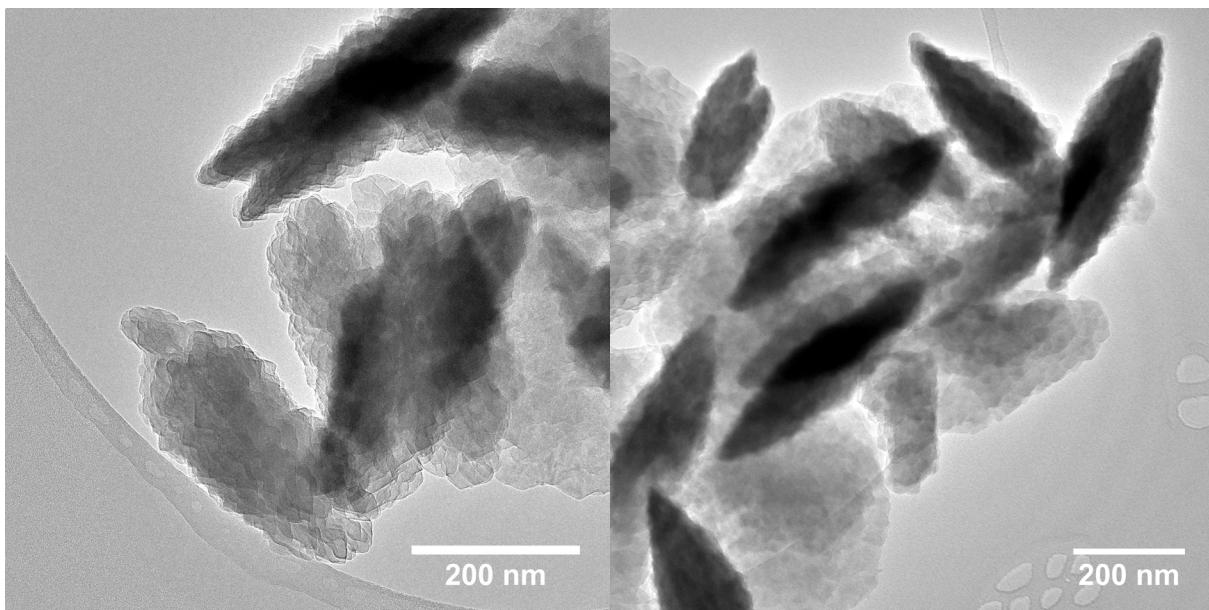


**Figure S2:** SEM image of CHA(2.3) zeolite

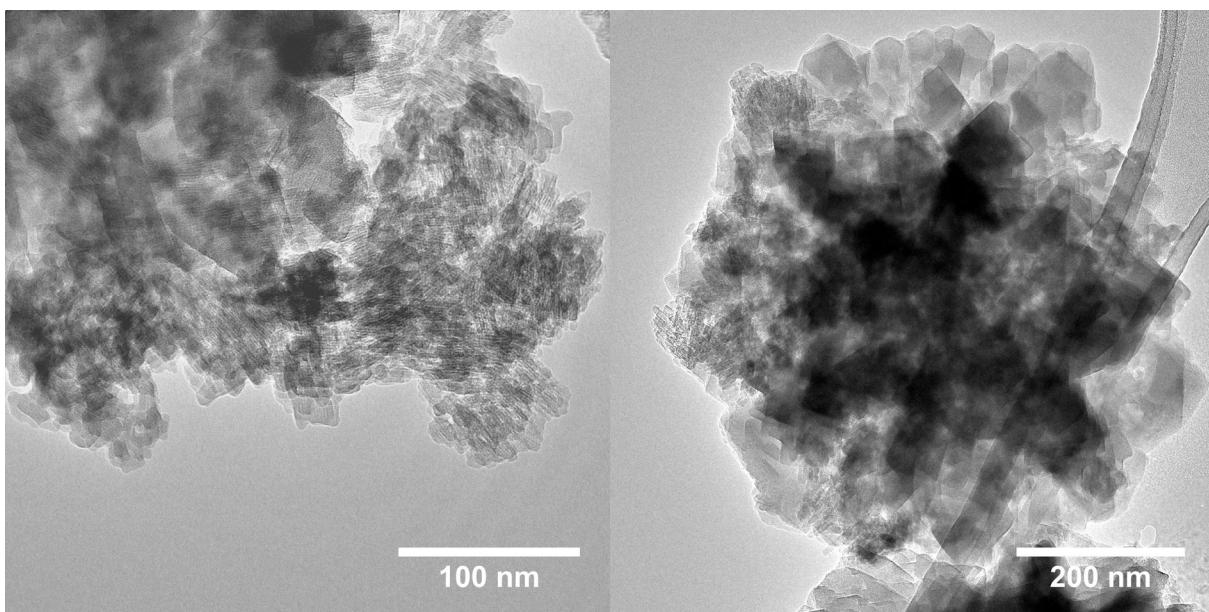


**Figure S3:** SEM image of CHA(1.9) zeolite

### TEM images

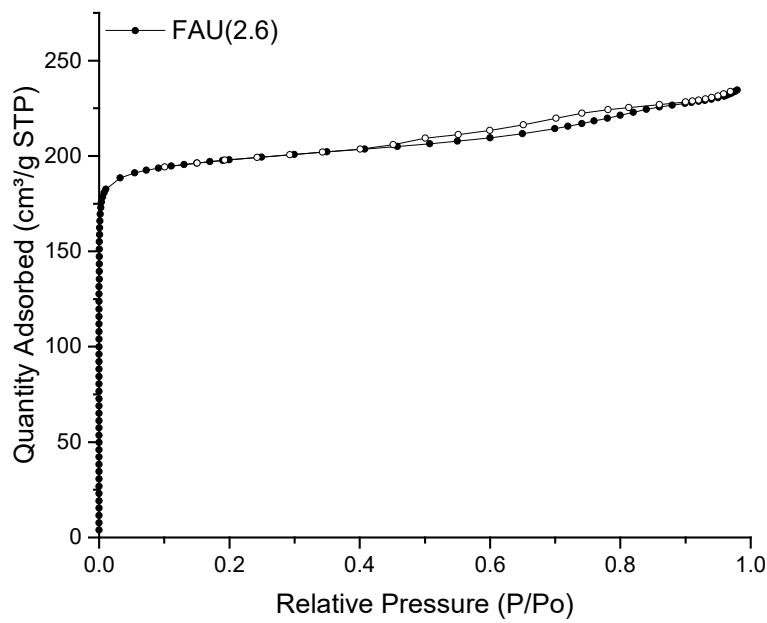


**Figure S4:** TEM image of CHA(1.9)

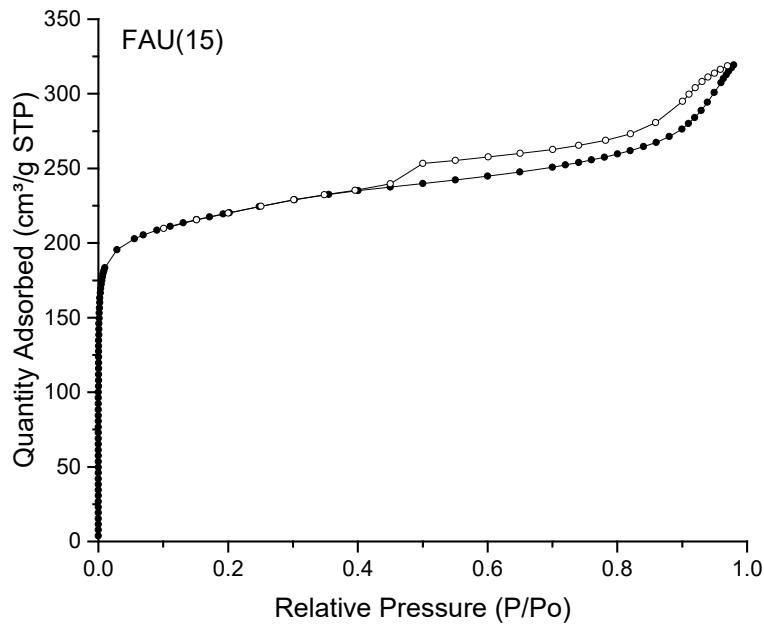


**Figure S5:** TEM image of CHA(2.3)

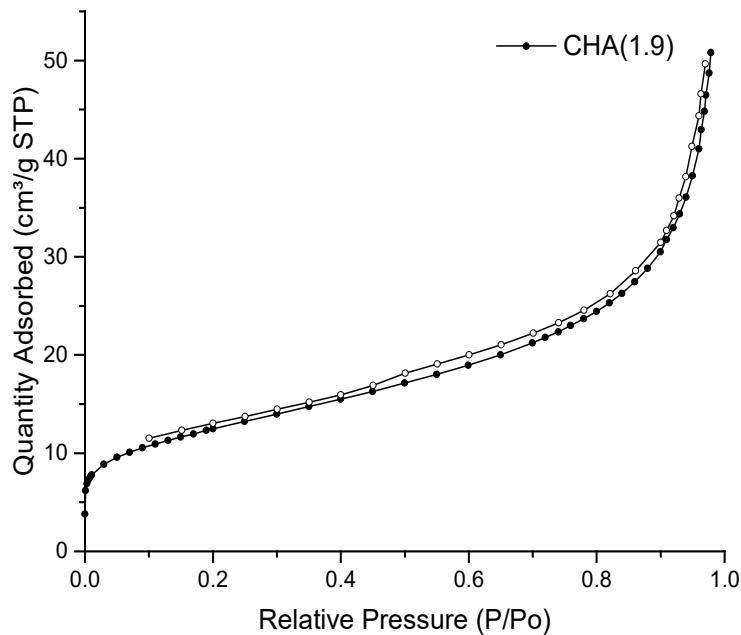
**Nitrogen physisorption of FAU(2.6), FAU(15), CHA(1.9), and CHA(2.3) zeolite samples**



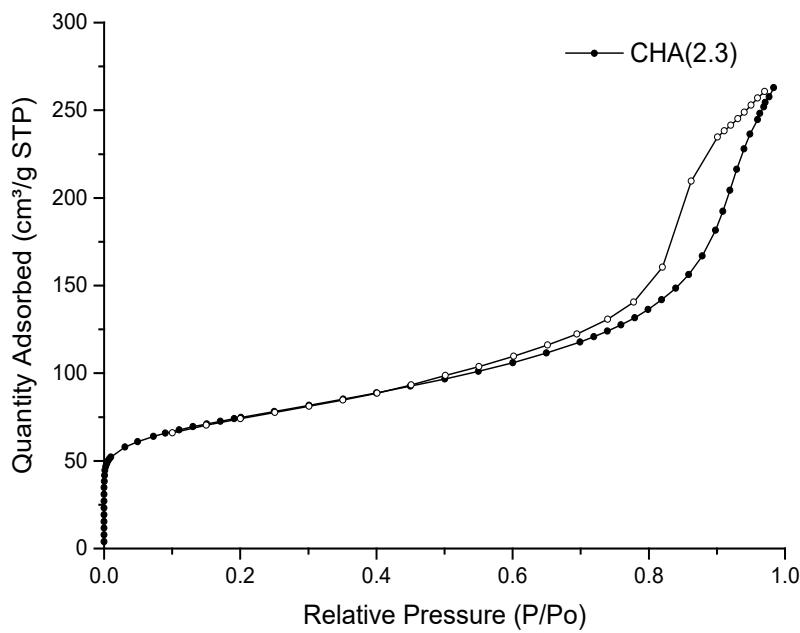
**Figure S6:** Nitrogen physisorption isotherm of FAU(2.6) zeolite at 77K



**Figure S7:** Nitrogen physisorption isotherm of FAU(15) zeolite at 77K



**Figure 8:** Nitrogen physisorption of CHA(1.9) at 77 K



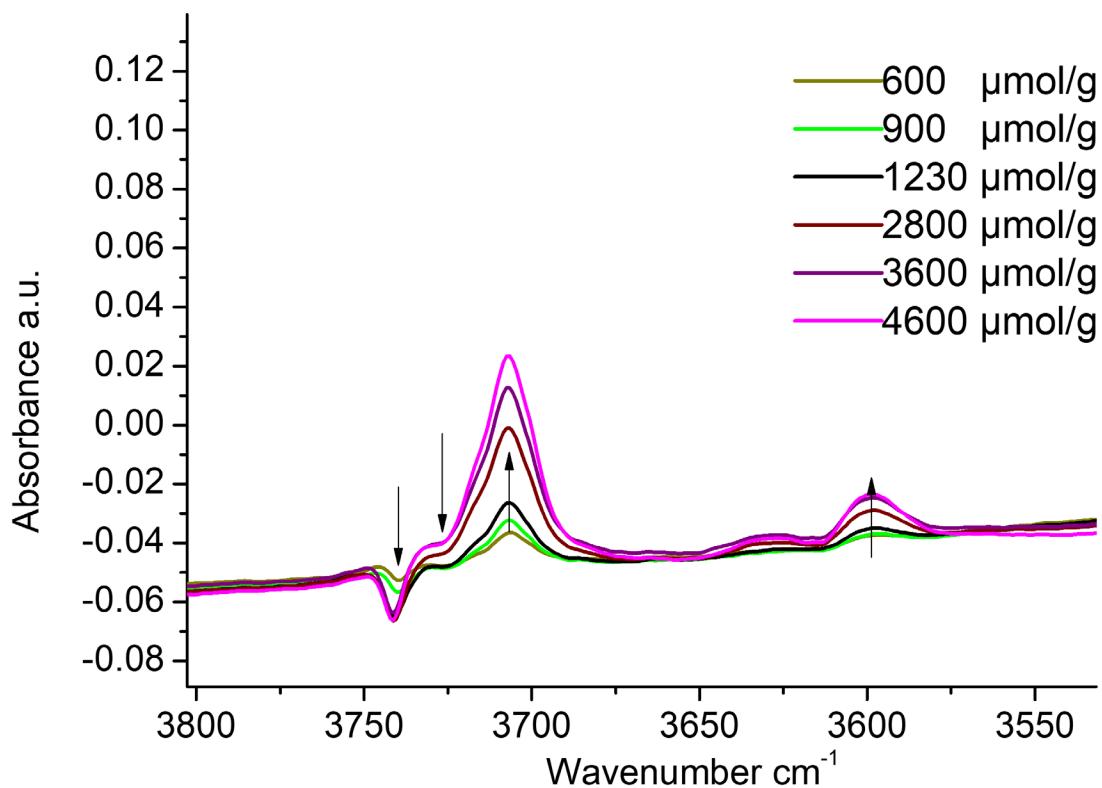
**Figure S9:** Nitrogen physisorption of CHA(2.3) at 77 K

**Table S3. Calculated porosity of FAU and CHA samples by N<sub>2</sub> physisorption at 77 K**

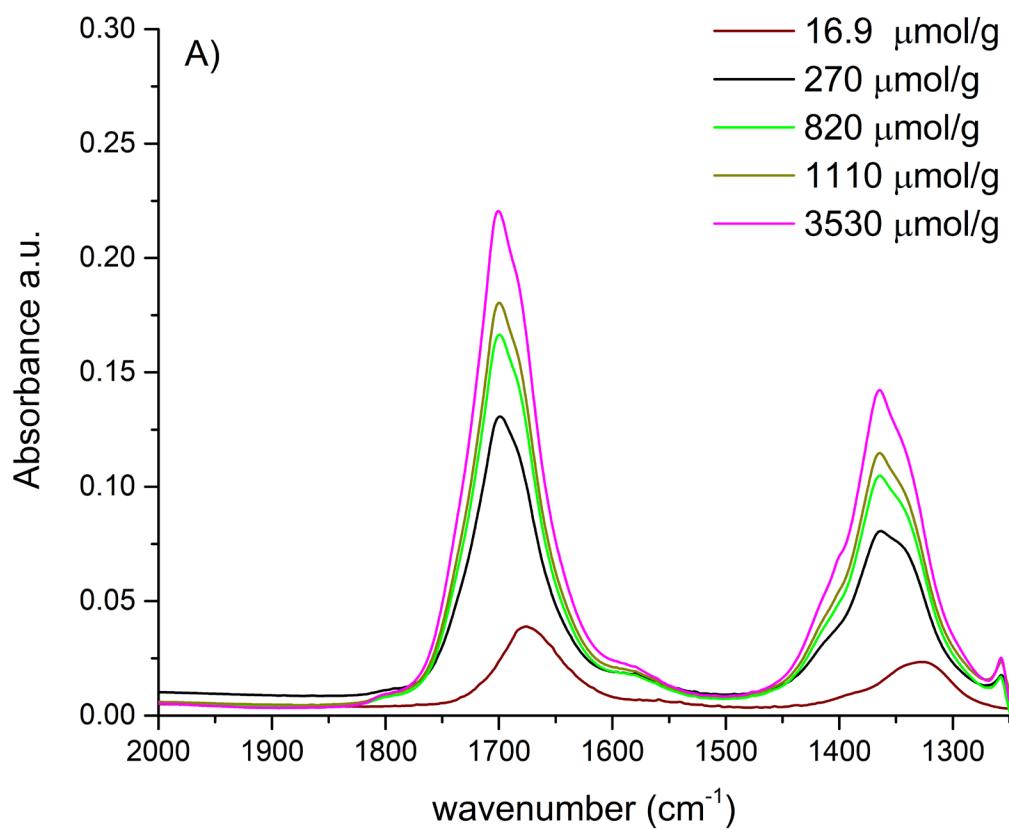
Sample	V <sub>total</sub> (cm <sup>3</sup> /g) <sup>a</sup>	V <sub>micropore</sub> (cm <sup>3</sup> /g) <sup>b</sup>	S <sub>BET</sub> (m <sup>2</sup> /g) <sup>c</sup>	S <sub>ext</sub> (m <sup>2</sup> /g) <sup>b,c</sup>
FAU(2.6)	0.49	0.280	695	140
FAU(15)	0.36	0.280	606	62
CHA(1.9)	0.07	0.004	43	33
CHA(2.3)	0.39	0.045	251	162

<sup>a</sup> Single point adsorption <sup>b</sup> Determined by the t-plot method <sup>c</sup> Determined by the Brunauer-Emmett-Teller method

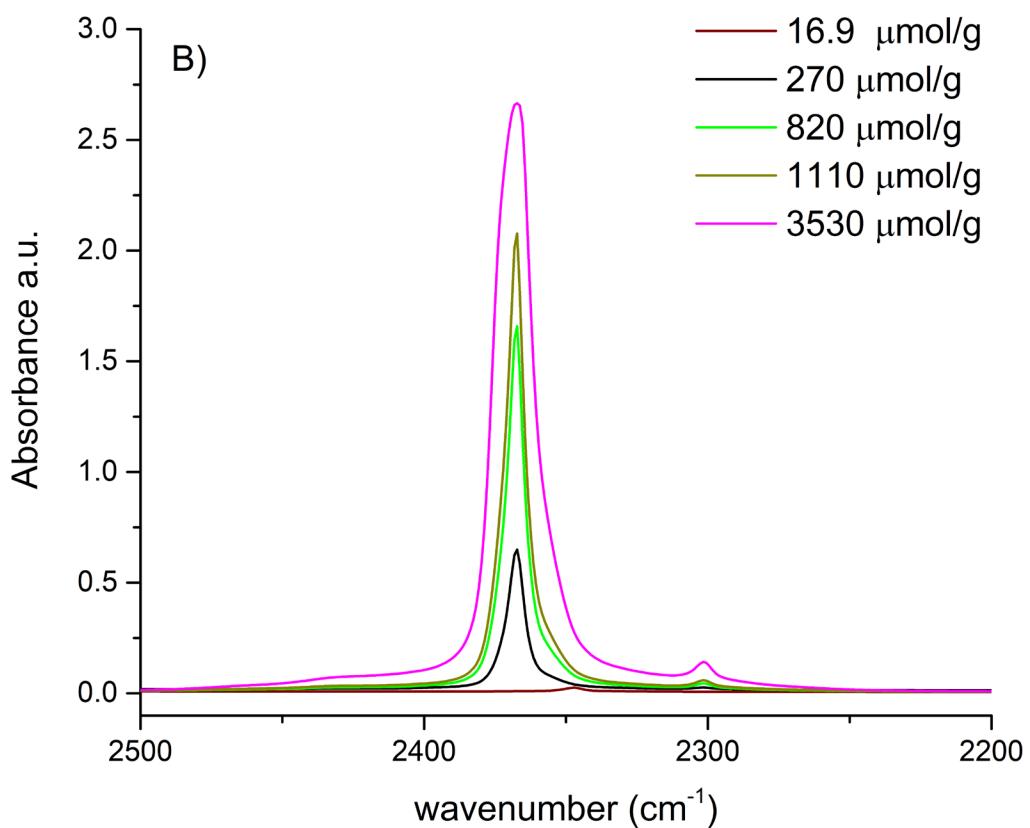
### CO<sub>2</sub> FTIR spectroscopic study



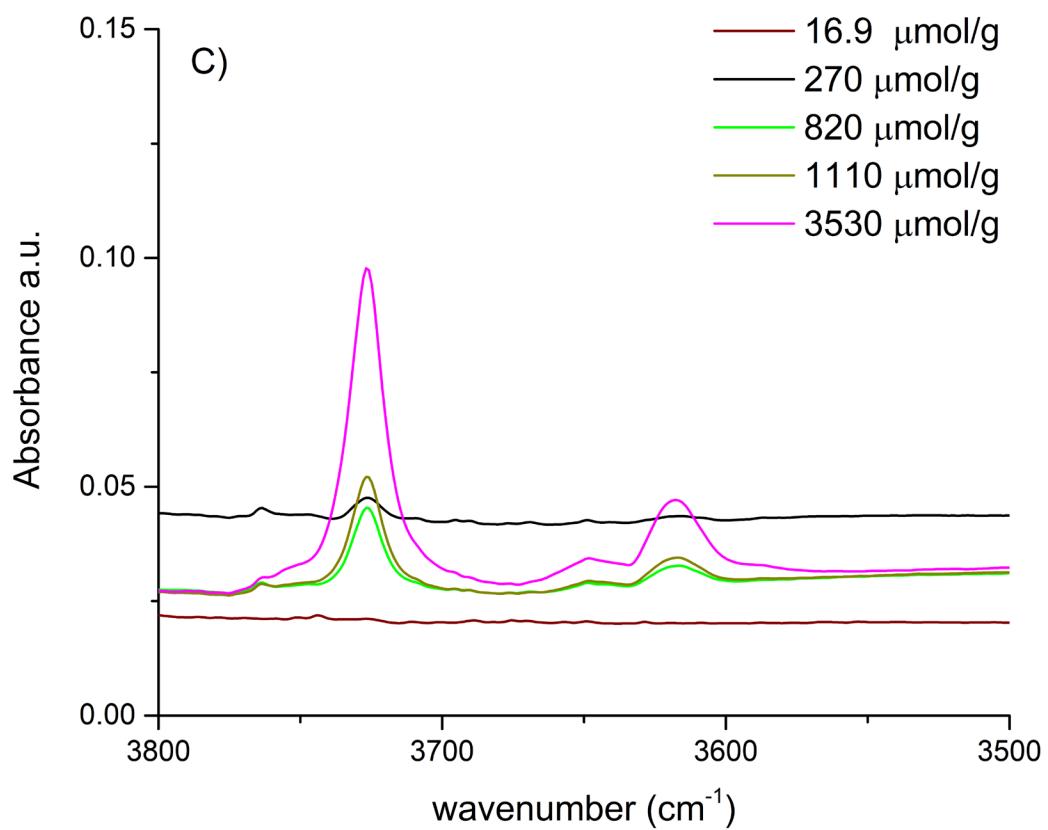
**Figure S10. IR spectra of CO<sub>2</sub> adsorbed on CHA(2.3) in the region of 3820-3510 cm<sup>-1</sup>**



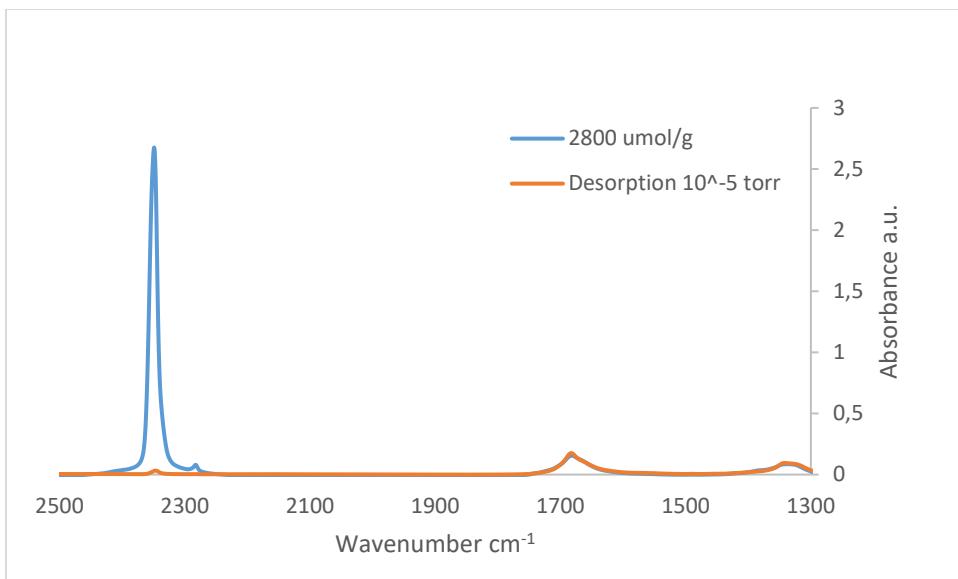
**Figure S11. IR spectra of pulsed CO<sub>2</sub> adsorbed on CHA(1.9) in the region of 2000-1250 cm<sup>-1</sup>**



**Figure S12. IR spectra of pulsed CO<sub>2</sub> adsorbed on CHA(1.9) in the region of 2500-2200 cm<sup>-1</sup>**



**Figure S13. IR spectra of pulsed CO<sub>2</sub> adsorbed on CHA(1.9) in the region of 3800-3500 cm<sup>-1</sup>**



**Figure S14: IR spectra of pulsed CO<sub>2</sub> adsorbed on CHA(2.3) and after desorption at 10<sup>-5</sup> torr**

## CO FTIR spectroscopic study

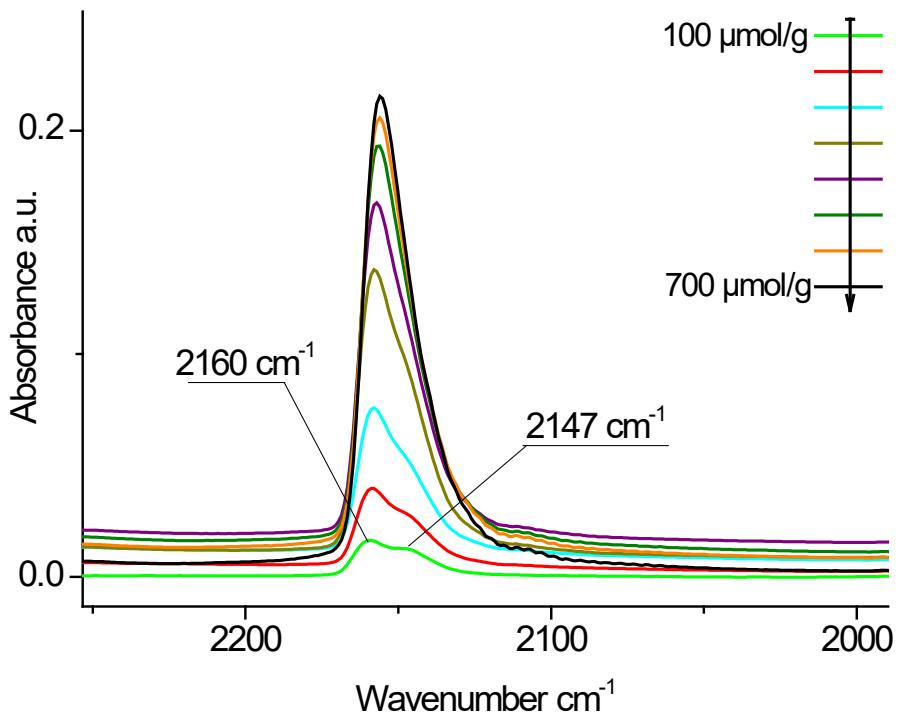


Figure S15. IR spectra of CO adsorbed on CHA(1.9) at 77 K.